Incrementalizing Random Sketching for Solving Consistent Linear Systems

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Outline

1. Overview: Linear System & Randomized Solvers
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OVERVIEW
Problem formulation

Our goal is to solve

\[
\begin{align*}
\begin{bmatrix} A \end{bmatrix}_{n \times d} & \begin{bmatrix} x \end{bmatrix}_d = \begin{bmatrix} b \end{bmatrix}_n, \\
\end{align*}
\]

for which we assume that at least one solution exists.

In particular, we are interested in randomized solvers, which recast the linear systems problem into a statistical estimation problem and then solve the estimation problem.
Two classes of randomized solvers

1. Random Sketching Methods. These methods use a matrix, $M$, with (much) fewer rows than $A$ and solve the problem

$$\min_x \| (MA)x - (Mb) \|^2_2,$$

where $M$ is a specifically structured random matrix, which we call a random sketching matrix.
2. Base Random Iteration. These methods use a random vectors, \( \{w_k : k + 1 \in \mathbb{N}\} \) and perform the iteration

\[
x_{k+1} = x_k + \gamma A'w_k [(w'_k b) - (w'_k A)x_k],
\]

where \( \gamma > 0 \) is some scalar.
Random sketching methods have the **lowest computational complexities** for finding an $\epsilon$-accurate solution (in residual) with high probability.
Why randomized solvers?

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Even with high-probability, we would think that the best solution then is the **random sketching** approach as it is the fastest method available.
The number of rows of $M$ depends on constants that are unknown and problem-specific, and is proportional to the reciprocal of $\epsilon$.

Therefore, $MA$, which can be cheap to compute (if $M$ is sparse), might still be too expensive to construct and store!
Challenges with Random Sketching

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TO SUMMARIZE

Random sketching has the best computational complexity, but we do not know how to choose the size of $M$ and it is nontrivial to store $MA$. 
OUR CONTRIBUTION
Overview & Consequences

We reformulate random sketching to implicitly construct $M A$ and simultaneously solve the projected system (i.e., $M A x = M b$). We do not need to decide on the size of $M$ a priori. We can let the size of $M$ grow implicitly, until, say, some stopping criteria is reached or the system is solved. Additionally, we do not need to create and store the matrix $M A$. We implicitly work with this matrix without constructing it.
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Additionally, we do not need to create and store the matrix $MA$. We implicitly work with this matrix without constructing it.
Owing to our reformulation, we are able to move towards the practical use of random sketching methods to solve actual linear systems.
OUR PROCEDURE
Step 1: Streaming rows of $M$

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Example: Gaussian Sketch. $M$ has independent, identically distributed standard Gaussian entries. Then $w_k$ is simply an $n$-dimensional standard Gaussian vector, and each $\{w_j\}$ are independent.
Step 2: Iterative Solver

Recalling that \( \{w_k\} \) are the rows of our sketching matrix \( M \), we now work through the iteration

\[
x_{k+1} = \begin{cases} 
    x_k + S_k A' w_k \frac{w_k' (b - A x_k)}{w_k' A S_k A' w_k} & \text{if } S_k A' w_k \neq 0 \\
    x_k & \text{otherwise}
\end{cases}
\]

and

\[
S_{k+1} = \begin{cases} 
    S_k - \frac{S_k A' w_k w_k' A S_k}{w_k' A S_k A' w_k} & \text{if } S_k A' w_k \neq 0 \\
    S_k & \text{otherwise}
\end{cases}
\]

where \( S_0 = I_d \) and \( x_0 \) is arbitrary.
Step 2: Iterative Solver

Note, \( \{S_k\} \) are orthogonal projections onto the space perpendicular to the rows of \( MA \) that have already been observed.
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In other words, \( R_I := \text{span}[A'w_0, \ldots, A'w_{l-1}] \) then \( S_l \) is an orthogonal projection onto \( R_I^\perp \).
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In other words, \( R_l := \text{span}[A'w_0, \ldots, A'w_{l-1}] \) then \( S_l \) is an orthogonal projection onto \( R_l^\perp \).

Therefore, as soon as we see the maximal possible linearly independent rows of \( MA \), then we will have solved the system to the highest accuracy possible allowed by \( M \) and \( A \).
THE CHALLENGE

How do we characterize this maximal set when the rows of $M$ are generated on the fly and they can have an arbitrary (independent, random permutation, adaptive, dependent) structure to previously observed rows of $M$?
THEORY
Subspace Characterizations

Let $\mathbf{w} \in \mathbb{R}^n$ be an arbitrary random variable. Define

$$\mathcal{N}(\mathbf{w}) = \text{span} \left[ \mathbf{z} \in \mathbb{R}^d : \mathbb{P}[\mathbf{z}' \mathbf{A}' \mathbf{w} = 0] = 1 \right]$$

(6)

and

$$\mathcal{R}(\mathbf{w}) = \mathcal{N}(\mathbf{w})^\perp.$$

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Lemma

$\mathcal{R}(\mathbf{w})$ is the smallest subspace of $\mathbb{R}^d$ such that $\mathbb{P}[\mathbf{A}' \mathbf{w} \in \mathcal{R}(\mathbf{w})] = 1.$
Subspace Characterizations

Message: For an arbitrary random variable $w$

- $\mathcal{R}(w)$ characterizes the row space of $w'A$
- $\mathcal{N}(w)$ characterizes the null space of $w'A$
- $\mathcal{V}(w)$ characterizes the deficiency of $w'A$ compared to $A$. 
Now for an arbitrary random variable $w$, let $\mathcal{R}(w)$ and $\mathcal{N}(w)$ be defined as before. For (a not necessarily related) set of random variables $\{w_k\}$, define

$$T = \min\{k \geq 0 : \text{span} [A'w_0, \ldots, A'w_k] \supset \mathcal{R}(w)\}.$$  

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$$T = \min\{k \geq 0 : \text{span} [A'w_0, \ldots, A'w_k] \supset \mathcal{R}(w)\}. \tag{8}$$

Again, we have not imposed any relationship between $w, w_0, w_1, \ldots$. Therefore, $T$ is quite generally defined. (This is useful when we consider parallel implementations.)
Theorem

Let $\mathbf{w}$ be a random variable, and let $\{\mathbf{w}_k\}$ be random variables such that $\mathbb{P}[A'\mathbf{w}_l \in \mathcal{R}(\mathbf{w})] = 1$ for all $l \geq 0$. Define $T$ as above. On the event $\{T < \infty\}$,

- For any $s \geq T + 1$, $S_{t+1} = S_s$ and $x_{T+1} = x_s$.
- If $A\mathbf{x} = \mathbf{b}$ admits a solution $\mathbf{x}^*$ (not necessarily unique), then

$$x_{T+1} = P_{\mathcal{N}(\mathbf{w})}\mathbf{x}_0 + P_{\mathcal{R}(\mathbf{w})}\mathbf{x}^*. \quad (9)$$
Do we solve the system?

**Corollary**

Under the settings of the preceding theorem, on the event \( \{ T < \infty \} \), \( Ax_{T+1} = b \) if and only if \( P_{\mathcal{V}(w)}x_0 = P_{\mathcal{V}(w)}x^* \).
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(1) When is \( T < \infty \)?

(2) When will \( P_{\mathcal{V}(w)}(x^* - x_0) = 0 \)?
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(1) When is \( T < \infty \)?

(2) When will \( P_{\mathcal{V}(w)}(x^* - x_0) = 0 \)?

Basically, when is this going to actually work?
When is this going to work?

Both of these questions will depend on how you choose $\mathbf{w}$, and how you design $\mathbf{w}_0$, $\mathbf{w}_1$, $\ldots$ for your particular system. This should depend on the linear system's structure and the hardware environment.
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Both of these questions will depend on how you choose $\mathbf{w}$, and how you design $\mathbf{w}_0, \mathbf{w}_1, \ldots$ for your particular system. This should depend on the linear system's structure and the hardware environment.

We have simply stated a very general theory of convergence for such methods, and supply specific examples in the paper.
SUMMARY

We restated matrix sketching as a random orthogonalization procedure and characterized the convergence for arbitrary sampling methodologies. This allows us to implicitly and incrementally generate and grow MA without storing it explicitly.